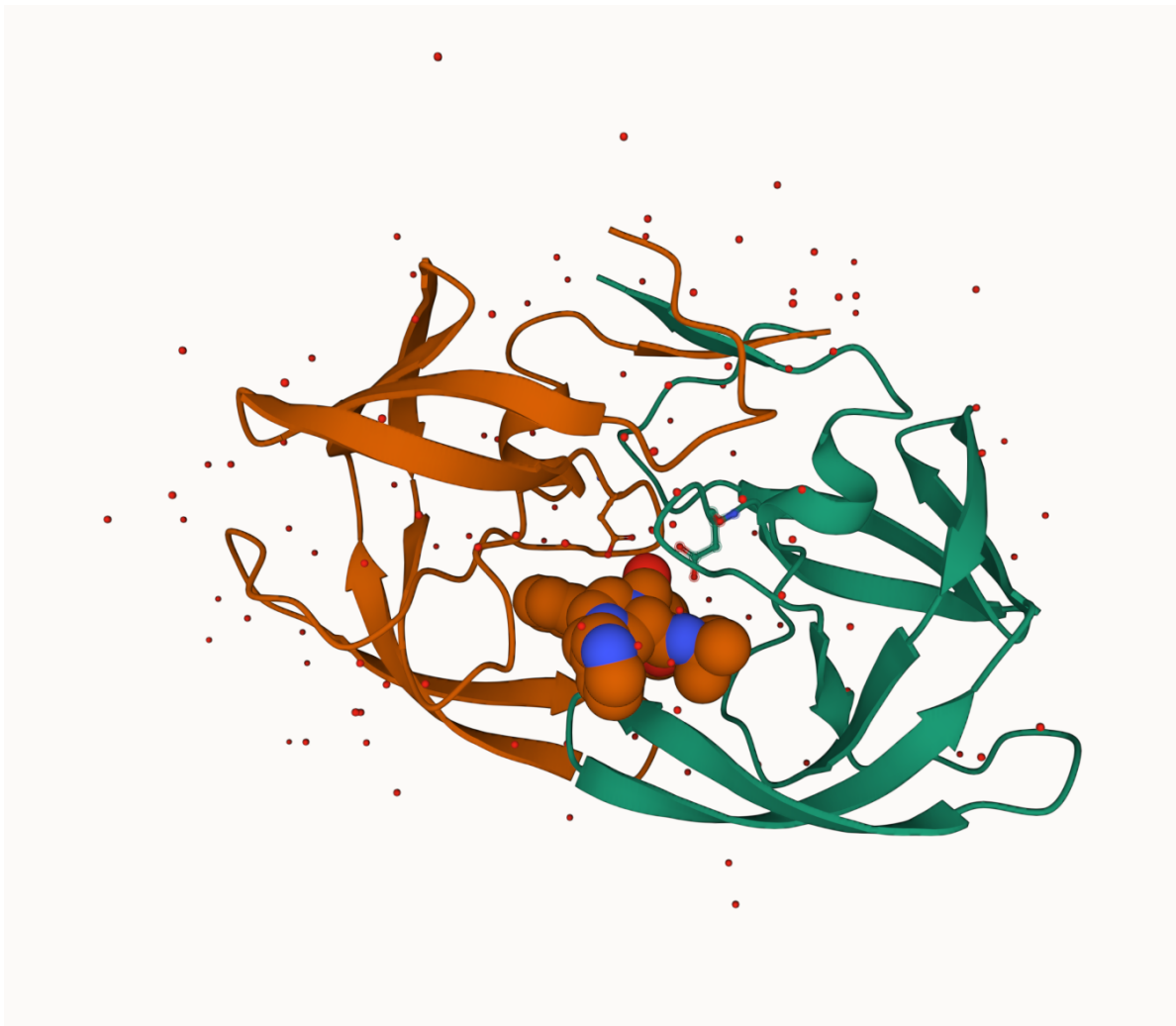


class9

Jie



```
library(bio3d)
```

```
pdb <- read.pdb("1hsg")
```

Note: Accessing on-line PDB file

```
pdb
```

```
Call: read.pdb(file = "1hsg")
```

```
Total Models#: 1
```

```
Total Atoms#: 1686, XYZs#: 5058 Chains#: 2 (values: A B)
```

```
Protein Atoms#: 1514 (residues/Calpha atoms#: 198)
```

```
Nucleic acid Atoms#: 0 (residues/phosphate atoms#: 0)
```

```
Non-protein/nucleic Atoms#: 172 (residues: 128)
```

```
Non-protein/nucleic resid values: [ HOH (127), MK1 (1) ]
```

```
Protein sequence:
```

```
PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYD  
QILIEICGHKAIGTVLVGPTPVNIIGRNLLTQIGCTLNFPQITLWQRPLVTIKIGGQLKE  
ALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYDQILIEICGHKAIGTVLVGPTP  
VNIIGRNLLTQIGCTLNF
```

```
+ attr: atom, xyz, seqres, helix, sheet,  
      calpha, remark, call
```

Q7: How many amino acid residues are there in this pdb object? 198

Q8: Name one of the two non-protein residues? HOH

Q9: How many protein chains are in this structure? 2

```
attributes(pdb)
```

```
$names
```

```
[1] "atom" "xyz" "seqres" "helix" "sheet" "calpha" "remark" "call"
```

```
$class
```

```
[1] "pdb" "sse"
```

Normal mode analysis (NMA) is a bioinformatics method for predicting functional motions.

```
adk <- read.pdb("6s36")
```

Note: Accessing on-line PDB file

PDB has ALT records, taking A only, rm.alt=TRUE

```
adk
```

```
Call: read.pdb(file = "6s36")
```

```
Total Models#: 1
```

```
Total Atoms#: 1898, XYZs#: 5694 Chains#: 1 (values: A)
```

```
Protein Atoms#: 1654 (residues/Calpha atoms#: 214)
```

```
Nucleic acid Atoms#: 0 (residues/phosphate atoms#: 0)
```

```
Non-protein/nucleic Atoms#: 244 (residues: 244)
```

```
Non-protein/nucleic resid values: [ CL (3), HOH (238), MG (2), NA (1) ]
```

```
Protein sequence:
```

```
MRIILLGAPGAGKGTQAQFIMEKYGIPQISTGDMRLRAAVKSGSELGKQAKDIMDAGKLV  
TDELVIALVKERIAQEDCRNGFLLDGFPRTIPQADAMKEAGINVDYVLEFDVPDELIVDKI  
VGRRVHAPSGRVYHVKFNPPKVEGKDDVTGEELTTRKDDQEETVRKRLVEYHQM  
TAPLIG  
YYSKEAEAGNTKYAKVDGTPVAEVRADLEKILG
```

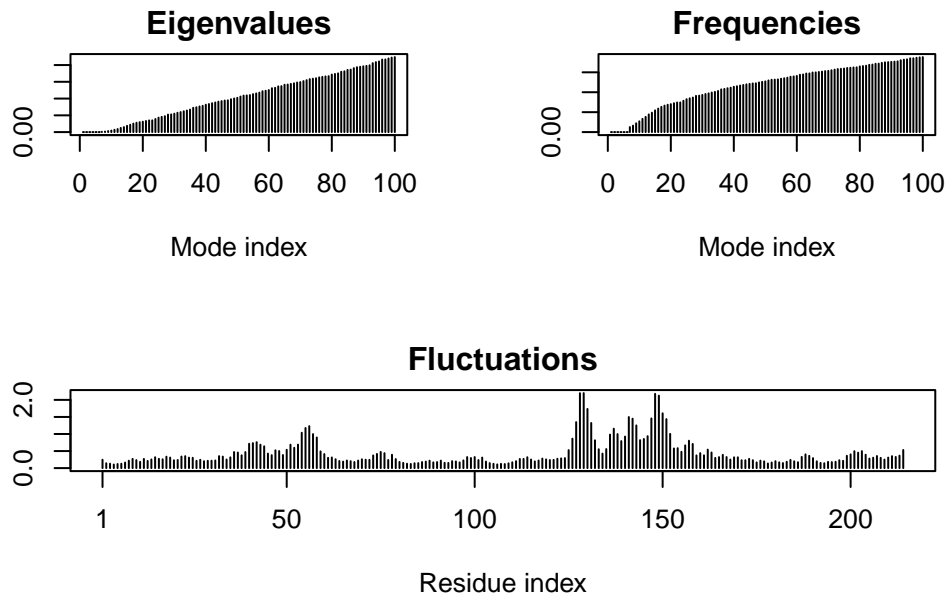
```
+ attr: atom, xyz, seqres, helix, sheet,  
      calpha, remark, call
```

```
m <- nma(adk)
```

```
Building Hessian... Done in 0.025 seconds.
```

```
Diagonalizing Hessian... Done in 0.834 seconds.
```

```
plot(m)
```



```
mktrj(m,file="adk_nma.pdb")
```

sequence of adk and use it to search the PDB database

```
aa <- get.seq("1ake_a")
```

Warning in get.seq("1ake_a"): Removing existing file: seqs.fasta

Fetching... Please wait. Done.

```
aa
```

```

      1      .      .      .      .      .      60
pdb|1AKE|A  MRIILLGAPGAGKGTQAQFIMEKYGIPQISTGDMRLRAAVKSGSELGKQAKDIMDAGKLV
      1      .      .      .      .      .      60

      61      .      .      .      .      .      120
pdb|1AKE|A  DELVIALVKERIAQEDCRNGFLLDGFPRTIPQADAMKEAGINVDYVLEFDVPDELIVDRI
      61      .      .      .      .      .      120

      121      .      .      .      .      .      180

```

```

pdb|1AKE|A    VGRRVHAPSGRVYHVKFNPPKVEGKDDVTGEELTTRKDDQEETVRKRLVEYHQMTAPLIG
              121          .          .          .          .          .          180

              181          .          .          .          214
pdb|1AKE|A    YYSKEAEAGNTKYAKVDGTPVAEVRADLEKILG
              181          .          .          .          214

```

Call:

```
read.fasta(file = outfile)
```

Class:

```
fasta
```

Alignment dimensions:

```
1 sequence rows; 214 position columns (214 non-gap, 0 gap)
```

+ attr: id, ali, call

Q13. How many amino acids are in this sequence, i.e. how long is this sequence? 214

```
#b <- blast.pdb(aa)
```

```
#hits <- plot(b)
```

```
hits <- NULL
```

```
hits$ pdb.id <- c('1AKE_A','6S36_A','6RZE_A','3HPR_A','1E4V_A','5EJE_A','1E4Y_A','3X2S_A',
```

```
pd.b.annotate(hits$ pdb.id)
```

	structureId	chainId	macromoleculeType	chainLength	experimentalTechnique
1AKE_A	1AKE	A	Protein	214	X-ray
6S36_A	6S36	A	Protein	214	X-ray
6RZE_A	6RZE	A	Protein	214	X-ray
3HPR_A	3HPR	A	Protein	214	X-ray
1E4V_A	1E4V	A	Protein	214	X-ray
5EJE_A	5EJE	A	Protein	214	X-ray
1E4Y_A	1E4Y	A	Protein	214	X-ray
3X2S_A	3X2S	A	Protein	214	X-ray
6HAP_A	6HAP	A	Protein	214	X-ray
6HAM_A	6HAM	A	Protein	214	X-ray

4K46_A	4K46	A	Protein	214	X-ray
3GMT_A	3GMT	A	Protein	230	X-ray
4PZL_A	4PZL	A	Protein	242	X-ray
	resolution	scopDomain			pfam
1AKE_A	2.00	Adenylate kinase	Adenylate kinase, active site lid (ADK_lid)		
6S36_A	1.60	<NA>	Adenylate kinase, active site lid (ADK_lid)		
6RZE_A	1.69	<NA>	Adenylate kinase, active site lid (ADK_lid)		
3HPR_A	2.00	<NA>	Adenylate kinase, active site lid (ADK_lid)		
1E4V_A	1.85	Adenylate kinase	Adenylate kinase, active site lid (ADK_lid)		
5EJE_A	1.90	<NA>	Adenylate kinase, active site lid (ADK_lid)		
1E4Y_A	1.85	Adenylate kinase	Adenylate kinase, active site lid (ADK_lid)		
3X2S_A	2.80	<NA>	Adenylate kinase, active site lid (ADK_lid)		
6HAP_A	2.70	<NA>	Adenylate kinase, active site lid (ADK_lid)		
6HAM_A	2.55	<NA>	Adenylate kinase, active site lid (ADK_lid)		
4K46_A	2.01	<NA>	Adenylate kinase, active site lid (ADK_lid)		
3GMT_A	2.10	<NA>	Adenylate kinase, active site lid (ADK_lid)		
4PZL_A	2.10	<NA>	Adenylate kinase, active site lid (ADK_lid)		

	ligandId
1AKE_A	AP5
6S36_A	CL (3),NA,MG (2)
6RZE_A	NA (3),CL (2)
3HPR_A	AP5
1E4V_A	AP5
5EJE_A	AP5,CO
1E4Y_A	AP5
3X2S_A	JPY (2),AP5,MG
6HAP_A	AP5
6HAM_A	AP5
4K46_A	ADP,AMP,PO4
3GMT_A	SO4 (2)
4PZL_A	CA,FMT,GOL

	ligandName
1AKE_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE
6S36_A	CHLORIDE ION (3),SODIUM ION,MAGNESIUM ION (2)
6RZE_A	SODIUM ION (3),CHLORIDE ION (2)
3HPR_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE
1E4V_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE
5EJE_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE,COBALT (II) ION
1E4Y_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE
3X2S_A	N-(pyren-1-ylmethyl)acetamide (2),BIS(ADENOSINE)-5'-PENTAPHOSPHATE,MAGNESIUM ION
6HAP_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE
6HAM_A	BIS(ADENOSINE)-5'-PENTAPHOSPHATE
4K46_A	ADENOSINE-5'-DIPHOSPHATE,ADENOSINE MONOPHOSPHATE,PHOSPHATE ION

3GMT_A		SULFATE ION (2)
4PZL_A		CALCIUM ION,FORMIC ACID,GLYCEROL
	source	
1AKE_A	Escherichia coli	
6S36_A	Escherichia coli	
6RZE_A	Escherichia coli	
3HPR_A	Escherichia coli K-12	
1E4V_A	Escherichia coli	
5EJE_A	Escherichia coli 0139:H28 str. E24377A	
1E4Y_A	Escherichia coli	
3X2S_A	Escherichia coli str. K-12 substr. MDS42	
6HAP_A	Escherichia coli 0139:H28 str. E24377A	
6HAM_A	Escherichia coli K-12	
4K46_A	Photobacterium profundum	
3GMT_A	Burkholderia pseudomallei 1710b	
4PZL_A	Francisella tularensis subsp. tularensis SCHU S4	

1AKE_A	STRUCTURE OF THE COMPLEX BETWEEN ADENYLATE KINASE FROM ESCHERICHIA COLI AND THE INHIB
6S36_A	
6RZE_A	
3HPR_A	
1E4V_A	
5EJE_A	
1E4Y_A	
3X2S_A	
6HAP_A	
6HAM_A	
4K46_A	
3GMT_A	
4PZL_A	

		citation	rObserved	rFree
1AKE_A	Muller, C.W., et al. J Mol Biol (1992)	0.19600	NA	
6S36_A	Rogne, P., et al. Biochemistry (2019)	0.16320	0.23560	
6RZE_A	Rogne, P., et al. Biochemistry (2019)	0.18650	0.23500	
3HPR_A	Schrank, T.P., et al. Proc Natl Acad Sci U S A (2009)	0.21000	0.24320	
1E4V_A	Muller, C.W., et al. Proteins (1993)	0.19600	NA	
5EJE_A	Kovermann, M., et al. Proc Natl Acad Sci U S A (2017)	0.18890	0.23580	
1E4Y_A	Muller, C.W., et al. Proteins (1993)	0.17800	NA	
3X2S_A	Fujii, A., et al. Bioconjug Chem (2015)	0.20700	0.25600	
6HAP_A	Kantaev, R., et al. J Phys Chem B (2018)	0.22630	0.27760	
6HAM_A	Kantaev, R., et al. J Phys Chem B (2018)	0.20511	0.24325	
4K46_A	Cho, Y.-J., et al. To be published	0.17000	0.22290	
3GMT_A	Buchko, G.W., et al. Biochem Biophys Res Commun (2010)	0.23800	0.29500	

4PZL_A						Tan, K., et al. To be published	0.19360 0.23680
	rWork	spaceGroup					
1AKE_A	0.19600	P 21 2 21					
6S36_A	0.15940	C 1 2 1					
6RZE_A	0.18190	C 1 2 1					
3HPR_A	0.20620	P 21 21 2					
1E4V_A	0.19600	P 21 2 21					
5EJE_A	0.18630	P 21 2 21					
1E4Y_A	0.17800	P 1 21 1					
3X2S_A	0.20700	P 21 21 21					
6HAP_A	0.22370	I 2 2 2					
6HAM_A	0.20311	P 43					
4K46_A	0.16730	P 21 21 21					
3GMT_A	0.23500	P 1 21 1					
4PZL_A	0.19130	P 32					

```
files <- get.pdb(hits$ pdb.id, path="pdbs", split=TRUE, gzip=TRUE)
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
1AKE.pdb.gz exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdb", split = TRUE, gzip = TRUE): pdbc/
6S36.pdb.gz exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdb", split = TRUE, gzip = TRUE): pdbs/
6RZE.pdb.gz exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
3HPR.pdb.gz exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdb", split = TRUE, gzip = TRUE): pdbs/
1E4V.pdb.gz exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdb", split = TRUE, gzip = TRUE): pdbs/
5EJE.pdb.gz exists. Skipping download
```

```
Warning in get.pdb(hits$pdb.id, path = "pdb", split = TRUE, gzip = TRUE): pdbs/
1E4Y.pdb.gz exists. Skipping download
```


Warning in get.pdb(hits\$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
3X2S.pdb.gz exists. Skipping download

Warning in get.pdb(hits\$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
6HAP.pdb.gz exists. Skipping download

Warning in get.pdb(hits\$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
6HAM.pdb.gz exists. Skipping download

Warning in get.pdb(hits\$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
4K46.pdb.gz exists. Skipping download

Warning in get.pdb(hits\$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
3GMT.pdb.gz exists. Skipping download

Warning in get.pdb(hits\$pdb.id, path = "pdbs", split = TRUE, gzip = TRUE): pdbs/
4PZL.pdb.gz exists. Skipping download

	0%
=====	8%
=====	15%
=====	23%
=====	31%
=====	38%
=====	46%
=====	54%
=====	62%
=====	69%
=====	77%

```
|
|=====| 85%
|
|=====| 92%
|
|=====| 100%
```

```
pdbbs <- pdbaln(files,fit=TRUE,exefile="msa")
```

Reading PDB files:

```
pdbbs/split_chain/1AKE_A.pdb
pdbbs/split_chain/6S36_A.pdb
pdbbs/split_chain/6RZE_A.pdb
pdbbs/split_chain/3HPR_A.pdb
pdbbs/split_chain/1E4V_A.pdb
pdbbs/split_chain/5EJE_A.pdb
pdbbs/split_chain/1E4Y_A.pdb
pdbbs/split_chain/3X2S_A.pdb
pdbbs/split_chain/6HAP_A.pdb
pdbbs/split_chain/6HAM_A.pdb
pdbbs/split_chain/4K46_A.pdb
pdbbs/split_chain/3GMT_A.pdb
pdbbs/split_chain/4PZL_A.pdb
  PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
..  PDB has ALT records, taking A only, rm.alt=TRUE
.... PDB has ALT records, taking A only, rm.alt=TRUE
.   PDB has ALT records, taking A only, rm.alt=TRUE
...
```

Extracting sequences

```
pdb/seq: 1   name: pdbbs/split_chain/1AKE_A.pdb
  PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 2   name: pdbbs/split_chain/6S36_A.pdb
  PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 3   name: pdbbs/split_chain/6RZE_A.pdb
  PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 4   name: pdbbs/split_chain/3HPR_A.pdb
```

```

PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 5   name: pdbc/split_chain/1E4V_A.pdb
pdb/seq: 6   name: pdbc/split_chain/5EJE_A.pdb
PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 7   name: pdbc/split_chain/1E4Y_A.pdb
pdb/seq: 8   name: pdbc/split_chain/3X2S_A.pdb
pdb/seq: 9   name: pdbc/split_chain/6HAP_A.pdb
pdb/seq: 10  name: pdbc/split_chain/6HAM_A.pdb
PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 11  name: pdbc/split_chain/4K46_A.pdb
PDB has ALT records, taking A only, rm.alt=TRUE
pdb/seq: 12  name: pdbc/split_chain/3GMT_A.pdb
pdb/seq: 13  name: pdbc/split_chain/4PZL_A.pdb

```

pdbc

```

[Truncated_Name:1] 1AKE_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:2] 6S36_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:3] 6RZE_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:4] 3HPR_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:5] 1E4V_A.pdb      1          .          .          .          40
-----MRIILLGAPVAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:6] 5EJE_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:7] 1E4Y_A.pdb      1          .          .          .          40
-----MRIILLGALVAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:8] 3X2S_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:9] 6HAP_A.pdb      1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:10] 6HAM_A.pdb     1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMEKYGIPQIS
[Truncated_Name:11] 4K46_A.pdb     1          .          .          .          40
-----MRIILLGAPGAGKGTQAQFIMAKFGIPQIS
[Truncated_Name:12] 3GMT_A.pdb     1          .          .          .          40
-----MRLILLGAPGAGKGTQANFIKEKFGIPQIS
[Truncated_Name:13] 4PZL_A.pdb     1          .          .          .          40
-----TENLYFQSNAMRIILLGAPGAGKGTQAKIIEQKYNIAHIS
                **^*****  *****  *  *^ *  **
41          .          .          .          80
TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDACKLVDELVIALVKE
TGDMLRAAVKSGSELGKQAKDIMDAGKLVDELVIALVKE

```

[Truncated_Name:8] 3X2S_A.pdb	TGDMRLRAAVKSGSELGKQAKDIMDCGKLVTDELVIALVKE
[Truncated_Name:9] 6HAP_A.pdb	TGDMRLRAAVKSGSELGKQAKDIMDAGKLVTDELVIALVRE
[Truncated_Name:10] 6HAM_A.pdb	TGDMRLRAAIKSGSELGKQAKDIMDAGKLVTDEIIIALVKE
[Truncated_Name:11] 4K46_A.pdb	TGDMRLRAAIKAGTELGKQAKSVIDAGQLVSDDIILGLVKE
[Truncated_Name:12] 3GMT_A.pdb	TGDMRLRAAVKAGTPLGVEAKTYMDEGKLVPDSLIIIGLVKE
[Truncated_Name:13] 4PZL_A.pdb	TGDMIRETIKSGSALGQELKKVLDAGELVSDEFIIKIVKD
	****~* ~* *~ ** * ~* ** * ~~ ~*~~
	41 . . . 80
	81 . . . 120
[Truncated_Name:1] 1AKE_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:2] 6S36_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:3] 6RZE_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:4] 3HPR_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:5] 1E4V_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:6] 5EJE_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:7] 1E4Y_A.pdb	RIAQEDCRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:8] 3X2S_A.pdb	RIAQEDSRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:9] 6HAP_A.pdb	RICQEDSRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:10] 6HAM_A.pdb	RICQEDSRNGFLLDGFPR TIPQADAMKEAGINVDYVLEFD
[Truncated_Name:11] 4K46_A.pdb	RIAQDDCAKGFLLDGFPR TIPQADGLKEVGVVVDYVIEFD
[Truncated_Name:12] 3GMT_A.pdb	RLKEADCANGYLFDFPR TIPQADAMKEAGVAIDYVLEID
[Truncated_Name:13] 4PZL_A.pdb	RISKNDCNNGFLLDGVPR TIPQAQELDKLGVNIDYIVEVD
	*~ * ~* ** **** ** ^ ~* ~**~* *
	81 . . . 120
	121 . . . 160
[Truncated_Name:1] 1AKE_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:2] 6S36_A.pdb	VPDELIVDKIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:3] 6RZE_A.pdb	VPDELIVDAIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:4] 3HPR_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDGTG
[Truncated_Name:5] 1E4V_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:6] 5EJE_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:7] 1E4Y_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:8] 3X2S_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:9] 6HAP_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:10] 6HAM_A.pdb	VPDELIVDRIVGRRVHAPSGRVYHV KFNPPKVEGKDDVTG
[Truncated_Name:11] 4K46_A.pdb	VADSVIVERMAGRRRAHLASGR TYHNVYNPPKVEGKDDVTG
[Truncated_Name:12] 3GMT_A.pdb	VPFSEIIERMSGRRTHPASGR TYHV KFNPPKVEGKDDVTG
[Truncated_Name:13] 4PZL_A.pdb	VADNLLIERITGRRIH PASGR TYHTKFNPPKVADKDDVTG
	* ~~~ ^ *** * *** ** ^***** *** **
	121 . . . 160

```

161 . . . 200
[Truncated_Name:1] 1AKE_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:2] 6S36_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:3] 6RZE_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:4] 3HPR_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:5] 1E4V_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:6] 5EJE_A.pdb EELTTRKDDQEECVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:7] 1E4Y_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:8] 3X2S_A.pdb EELTTRKDDQEETVRKRLCEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:9] 6HAP_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:10] 6HAM_A.pdb EELTTRKDDQEETVRKRLVEYHQM TAPLIGYYSKEAEAGN
[Truncated_Name:11] 4K46_A.pdb EDLVIREDDKEETVLARLGVYHNQ TAPLIAYYGKEAEAGN
[Truncated_Name:12] 3GMT_A.pdb EPLVQRDDKEETVKKRLDVYEAQTK PLITYYGDWARRGA
[Truncated_Name:13] 4PZL_A.pdb EPLITRTDDNEDTVKQRLSVYHAQT AKLIDFYRNFSSNT
* * * * * ^ * * * * ^ *
161 . . . 200

```

```

201 . . 227
[Truncated_Name:1] 1AKE_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:2] 6S36_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:3] 6RZE_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:4] 3HPR_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:5] 1E4V_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:6] 5EJE_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:7] 1E4Y_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:8] 3X2S_A.pdb T--KYAKVDGTKPVAEVRADLEKILG-
[Truncated_Name:9] 6HAP_A.pdb T--KYAKVDGTKPVCEVRADLEKILG-
[Truncated_Name:10] 6HAM_A.pdb T--KYAKVDGTKPVCEVRADLEKILG-
[Truncated_Name:11] 4K46_A.pdb T--QYLKFDGTKAVAESAELEKALA-
[Truncated_Name:12] 3GMT_A.pdb E-----NGLKAPA-----YRKISG-
[Truncated_Name:13] 4PZL_A.pdb KIPKYIKINGDQAVEKVSQDIFDQLNK
*
201 . . 227

```

Call:

```
pdbaln(files = files, fit = TRUE, exefile = "msa")
```

Class:

```
pdbs, fasta
```

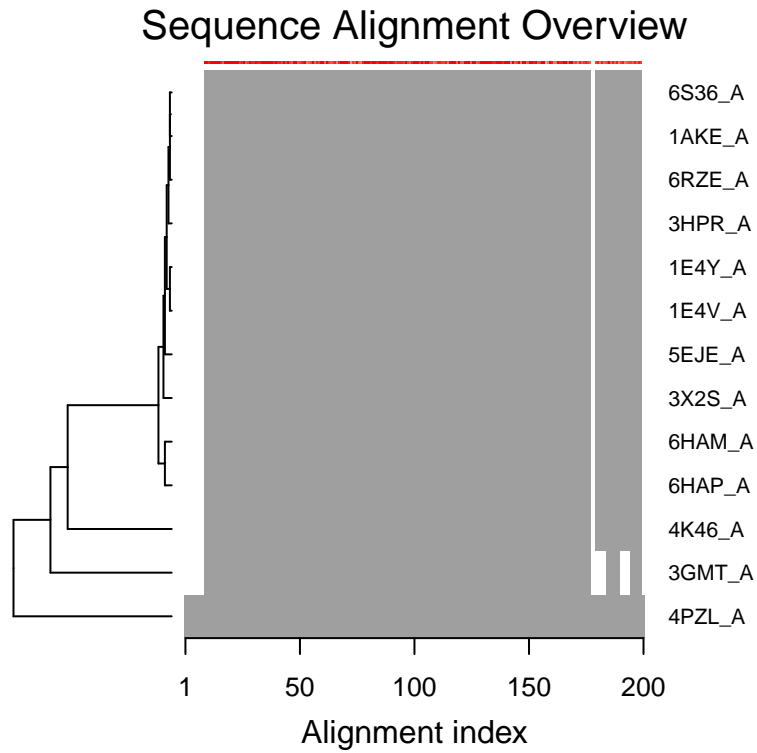
Alignment dimensions:

```
13 sequence rows; 227 position columns (204 non-gap, 23 gap)
```

```
+ attr: xyz, resno, b, chain, id, ali, resid, sse, call
```

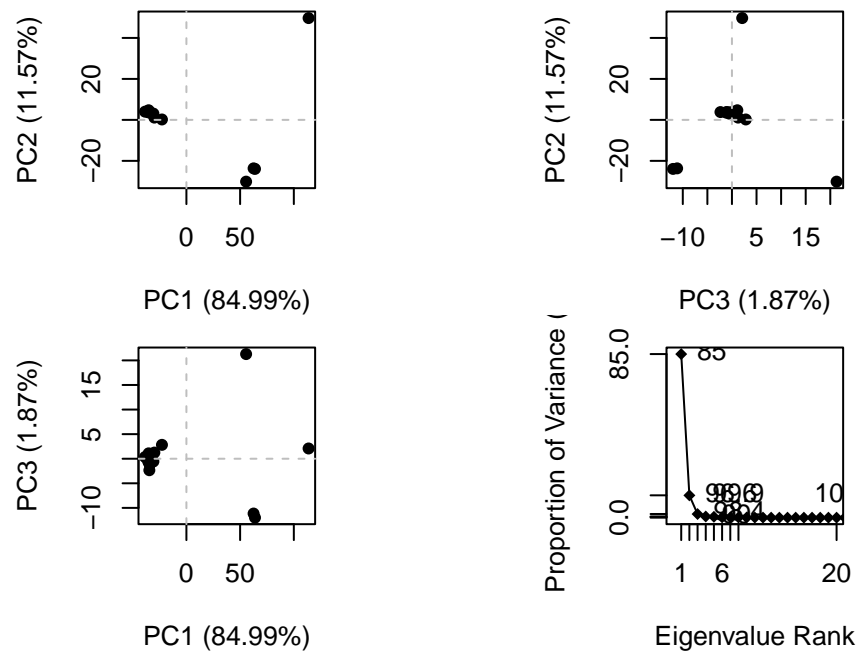
```
ids <- basename.pdb(pdb$id)
```

```
plot(pdb, labels=ids)
```



```
pc.xray <- pca(pdb)
```

```
plot(pc.xray)
```

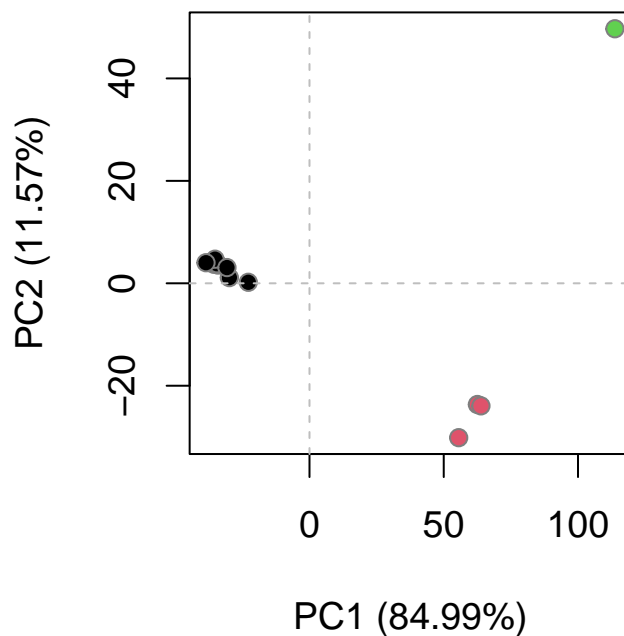


```
rd <- rmsd(pdbbs)
```

Warning in rmsd(pdbbs): No indices provided, using the 204 non NA positions

```
hc.rd <- hclust(dist(rd))
grps.rd <- cutree(hc.rd, k=3)

plot(pc.xray, 1:2, col="grey50", bg=grps.rd, pch=21, cex=1)
```



```
modes <- nma(pdb)
```

Details of Scheduled Calculation:

```
... 13 input structures
... storing 606 eigenvectors for each structure
... dimension of x$U.subspace: ( 612x606x13 )
... coordinate superposition prior to NM calculation
... aligned eigenvectors (gap containing positions removed)
... estimated memory usage of final 'eNMA' object: 36.9 Mb
```

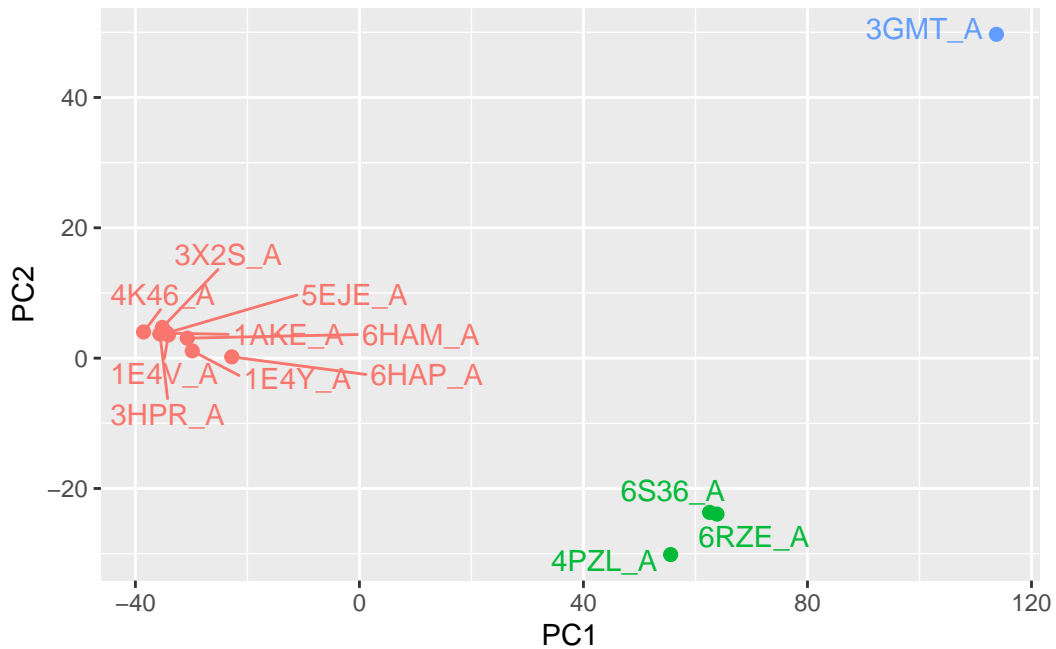
	0%
=====	8%
=====	15%
=====	23%
=====	31%



```
library(ggplot2)
library(ggrepel)

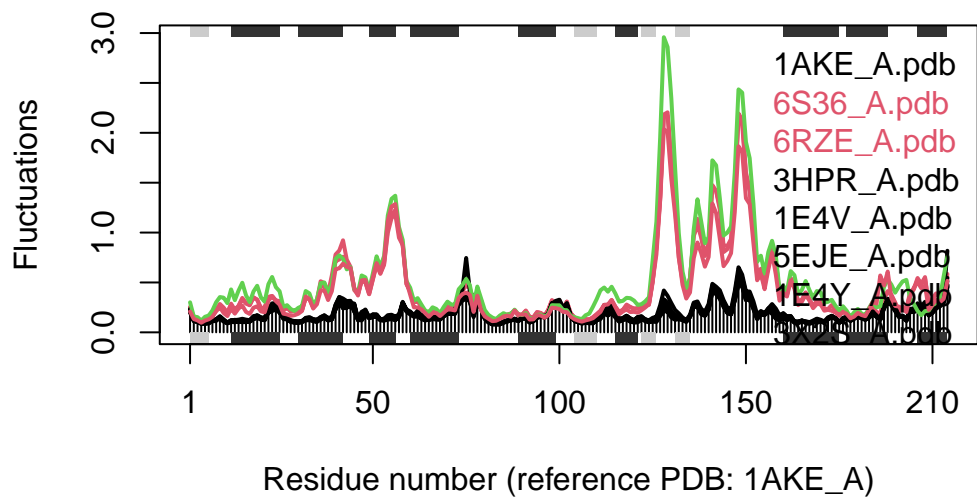
df <- data.frame(PC1=pc.xray$z[,1],
                  PC2=pc.xray$z[,2],
                  col=as.factor(grps.rd),
                  ids=ids)

p <- ggplot(df) +
  aes(PC1, PC2, col=col, label=ids) +
  geom_point(size=2) +
  geom_text_repel(max.overlaps = 20) +
  theme(legend.position = "none")
p
```



```
plot(modes, pdbs, col=grps.rd)
```

Extracting SSE from pdbs\$sse attribute



Q14. What do you note about this plot? Are the black and colored lines similar or different? Where do you think they differ most and why?

The colored lines (red and green) look similar but they are very different from the black lines. Colored lines also have higher and more peaks than the black lines. It indicates that there are two major distinct conformational states of Adenylate kinase. These differ by a collective low frequency displacement of two nucleotide-binding site regions.